

Technical breakdown of position based fluid

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This paper will implement a parallel version of Position Based Fluids (PBF) method [Miles Macklin et al. 2013] for real-time fluid simulation and discuss the effects of different confinement in the original paper. Compare to BPF, the widely used Lagrangian methods often sacrifice the incompressibility of fluids for better performance, which cause the result to be inaccurate compared to the behavior of real fluids. To address this issue, PBF method numerically integrates quantities from the neighboring particles with a smoothing kernel to calculate displacement while enforcing the incompressible constraint of fluid for the Navier-Stokes equations. This paper uses CUDA framework in our implementation to achieve real-time, accurate result.

CCS Concepts: • **Computer methodologies** → **Physical simulation**.

Additional Key Words and Phrases: Fluids, Physically Based Animation

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1 INTRODUCTION

Fluids such as water or gases, has played an important role in the field of computer based physical animations. Although fluid may look simple, but they are much more complicated to simulate than solids. The motion of fluids are governed by the Navier-Stokes equations. The solutions of the Navier-Stokes equations often include turbulence which makes finding an analytical solution extremely hard in three dimensions. Although countless efforts are put into this field, the progress is still limited. Currently numerical simulation remains the most effective way to approximate a complex Navier-Stokes system. In this report, we'll breakdown the Position based method and discuss the significance and effects of all the constraints it proposed. In position based fluid, fluids are divided into particles and then compute the interactions between the fluid particles such as density and viscosity confinement, combine with external forces, we then update the velocities and positions of the fluid particle. To optimize for better performance,

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we divided the simulation spaces into grids so that particles only interact with other particles in the neighboring grid.

2 RELATE WORK

According to simulation method, algorithm can generally be divided into two different categories: Euler and Lagrangian methods. Euler methods split the space into grids and use finite difference method to evaluate the model. One of the examples is Stam's stable fluid [Stam 1999]. The Lagrangian methods divide the fluid into free particles under constraints. One of the most known methods of this kind is the SPH method [Smoothed Particle Hydrodynamics]. Both categories have their own drawbacks: Euler methods are easily affected by numerical dissipation during simulation and Lagrangian methods need smaller steps each iteration to ensure the stability of the algorithm. Current fluid simulation methods such as Position Based Fluid methods use a hybrid of both methods to achieve a high level of accuracy while having reasonable performance.

3 PBF METHOD

For any arbitrary fluid, its motion is governed by the Navier-Stokes equation:

$$\rho \frac{\partial \mathbf{v}}{\partial t} = \rho \mathbf{g} - \nabla p + \mu \nabla^2 \mathbf{v} \quad (1)$$

where ρ is the density of the fluid, μ is the viscosity constant of the fluid, p is the pressure and \mathbf{v} denote the velocity. However, the above equation does not satisfy the incompressible property of fluids, so an additional constraint needs to be added:

$$\nabla \cdot \mathbf{v} = 0 \quad (2)$$

In PBF method, we represent the fluid with N particles, and their positions are denoted as $\mathbf{P} = \{\mathbf{p}_1, \mathbf{p}_2, \dots, \mathbf{p}_N\}$.

3.1 Density confinement

Since we are simulating a fluid, we have to satisfy the incompressible property of the fluid. Instead of satisfying equation (2) directly, BPF uses density constraint to satisfy the requirement:

$$C_i(\mathbf{p}_1, \mathbf{p}_2, \dots, \mathbf{p}_N) = \frac{\rho_i}{\rho_0} - 1 \quad (3)$$

Where ρ_0 is the rest density of the fluid and ρ_i is given by the standard SPH estimator:

$$\rho_i = \sum_j m_j W(\mathbf{p}_i - \mathbf{p}_j, h) \quad (4)$$

where m_j is the mass of the neighboring particles and W is the poly6 smoothing kernel as given in [Müller et al. 2003]. In this paper, we assume all the particles have a mass of 1. Since a fluid is incompressible, intuitively this means that

the density for any given part of the fluid should equal to the rest density of the fluid, which means the result of (3) should be equal to zero at all time:

$$C_i(\mathbf{p}_1, \mathbf{p}_2, \dots, \mathbf{p}_2) = 0 \quad (5)$$

Furthermore, when there is an displacement in the fluid particle, the density should also remains the same, this property is statisfied through the following constraint:

$$C(\mathbf{p} + \Delta\mathbf{p}) = 0 \quad (6)$$

We can estimate a proper $d\Delta\mathbf{p}$ equation (6) with Newton steps along the gradient of equation(5):

$$\Delta\mathbf{p} = \nabla C(\mathbf{p})\lambda \quad (7)$$

And we can approximate the value of equation (6) with linear approximation:

$$C(\mathbf{p} + \Delta\mathbf{p}) \approx C(\mathbf{p}) + \nabla C^T \nabla C \lambda \quad (8)$$

Now we only need the value of ∇C^T to estimate a proper λ . Luckily, ∇C^T can be defined as follows:

$$\nabla_{\mathbf{p}_k} C_i = \begin{cases} \sum_j \nabla_{\mathbf{p}_k} W(\mathbf{p}_i - \mathbf{p}_j, h) & \text{if } k = i \\ -\nabla_{\mathbf{p}_k} W(\mathbf{p}_i - \mathbf{p}_j, h) & \text{if } k = j \end{cases} \quad (9)$$

Plugging this back to (8) and solve for λ :

$$\lambda_i = \frac{\nabla_{\mathbf{p}_k} C_i}{\sum_k |\nabla_{\mathbf{p}_k} C_i|^2 + \epsilon} \quad (10)$$

Note that when there are not particles nearby, $\sum_k |\nabla_{\mathbf{p}_k} C_i|^2$ might be zero, so a extra ϵ term is added to prevent the denominator from evaluating to zero.

3.2 Tensile Instability

The original SPH algorithm has an issue known as particle clamping, where particles clamps to each other. Therefore, a corrective term s_{corr} is added when computing $\Delta\mathbf{p}$:

$$\Delta\mathbf{p}_i = \frac{1}{\rho_0} \sum_j (\lambda_i + \lambda_j + s_{corr}) \nabla W(\mathbf{p}_i - \mathbf{p}_j, h) \quad (11)$$

And it is define as follows:

$$s_{corr} = \left(\frac{-kW(\mathbf{p}_i - \mathbf{p}_j, h)}{W(\Delta\mathbf{q}, h)} \right)^n \quad (12)$$

Where k , n , h and $\Delta\mathbf{q}$ is selected constants. In our implementation, we use $k = 0.1$, $n = 4$, $h = 0.1$ and $|\Delta\mathbf{q}| = 0.1h$. With this corrective term, PBF does not required 30-40 neiboring particles at all time like SPH does, improcing simulation efficiency.

3.3 Vorticity and viscosity confinement

Since PBF introduced addtional damping which is undesirable, [Fedkiw et al. 2001] introduced vorticity confinement which has the following form:

$$\mathbf{f}_i^{\text{vorticity}} = \epsilon(\mathbf{N} \times \omega_i) \quad (13)$$

where

$$\omega_i = \sum_j (\mathbf{v}_j - \mathbf{v}_i) \times \nabla_{\mathbf{p}_j} W(\mathbf{p}_i - \mathbf{p}_j, h) \quad (14)$$

And $\mathbf{N} = \frac{\omega_i}{\nabla|\omega_i|}$.

Note that we did not implement this vorticity confinement in our code.

Compare to vorticity confinement, viscosity confinement is relatively simple. We apply the XSPH viscosity estimator to oue velocity update from [Schechter and Bridson 2012] to accomadate the coherent motion for any arbitrary fluid:

$$\mathbf{v}_i^{\text{new}} = \mathbf{v}_i + c \sum_j (\mathbf{v}_j - \mathbf{v}_i) W(\mathbf{p}_i - \mathbf{p}_j, h) \quad (15)$$

Equation (15) will introduce additional interation between neiboring particles so the movements of neiboring particle tend to "drag" the target particle to move with them.

4 RESULTS

We implemented the PBF method using CUDA framework. Note that in our implementation we omitted the vorticity confinement due to time limits. We simulated a sene where a block of water falls into a empty square space. Our simulation has 8000 particles in total and wer are able to render their movements in real time. The behavior of our simulated fluid agree with our expectation: The fluid particles having coherent motion while maintaing adquate distance with each other (less compression). Therefore we concluded that our simulation produces a accurate result.

5 DISCUSSION

In this paper, we covered the technical details of the PBF and provided a parallel implmentation using CUDA framework and a optimization inspired by [TODO: NVIDIA and tongji]. BPF method produces a realistic simulation to the human eye while maintaining a resonable performance. One limit of our implementation is that vorticity confinement is not implemented, which will cause the fluid to be more damped than in reality. In our implementation, we observed a performance drop of the simulation when the fluid particles start hitting the boundry. We conclude that this is due to particles concentrated in a few grid cells in space, hence more computation needed for each cells.

6 REFERENCES